

Figure 1: Association [3H]-4MG

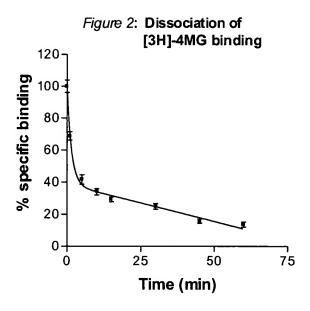


Figure 3: DRUG INHIBITION OF [3H]-4MG BINDING

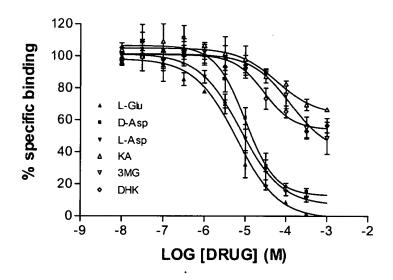
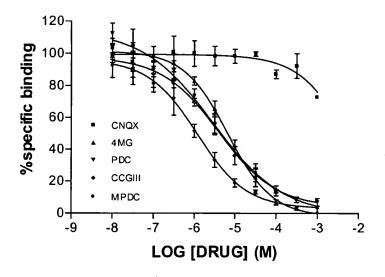


Figure 4: DRUG INHIBITION of [3H]-4MG BINDING



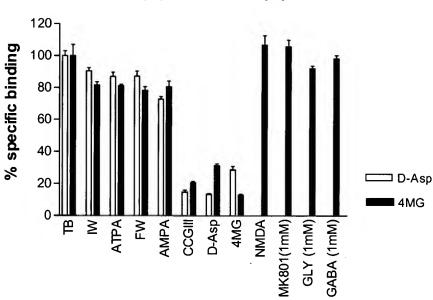


Figure 5: COM PARISON of MISCELLANEOUS DRUG INHIBITION of [3H]-D-ASPARTATE and [3H]-4MG

PhO₂SHNOC

$$O_2$$
H

 O_2 SHNOC

 O_2 H

 O_2 SHNOC

 O_2 H

 O_2 SHNOC

 O_2 H

 O_2 SHNOC

 O_2 H

 O_2 CONHSO₂CH₃
 O_2 SHNOC

 O_2 H

 O_2 CONHSO₂CH₃
 O_2 SHNOC

 O_2 H

 O_2 CONHSO₂CH₃
 O_2 SHNOC

 O_2 H

 O_2 CONHSO₂CH₃

$$Ph$$
 NH_2
 HO_2C
 CO_2H
 Ph
 NH_2
 HO_2C
 CO_2H

Figure 6A

Figure 6B

Figure 6C

CO2H

$$H_3CO_2C$$
 NH_2
 HO_2C
 CO_2H
 NH_2
 NH_2

NHCONHPh

$$OH$$
 OH
 OH

Figure 6D

Figure 6E

Figure 6F

Figure 6G

CO₂H

HO₂C



$$CO_2H$$
 NH_2
 NH_2

Figure 6H

$$RO_2C$$
 $R = Me, Et, ^{1}Bu$
 $R^{1}R^{2}$
 $R^{1}R^{2}$
 $R^{2}NR^{3}R^{4}$
 $R^{5}O_2C$
 $R^{5}O_2C$

 $R^1 = CH_3$, and halogen R^2 , R^3 are independently

H, C1-C6-alkyl, C3-C4-alkenyl, C3-C5-cycloalkyl, C1-C6-alkyl-CO-,

C1-C6-alkyl-OCO-, C1-C6-alkyl-NHCO-, HCO-, or C3-C6-alkynyl

 R^2 , R^3 taken together can be $-CH_2(CH_2)_pCH_2$ -

$$R$$
 CO_2H
 HO_2C
 $R = H, Me, Et, Cl$
 $R = H, Me, Et, nPr$
 $R = H, Et, nPr$